



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:03 PM GMT

PDB ID : 1WC5  
Title : SOLUBLE ADENYLYL CYCLASE CYAC FROM S. PLATENSIS IN COM-  
PLEX WITH ALPHA,BETA-METHYLENE-ATP IN PRESENCE OF  
BICARBONATE  
Authors : Steegborn, C.; Litvin, T.N.; Levin, L.R.; Buck, J.; Wu, H.  
Deposited on : 2004-11-08  
Resolution : 2.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

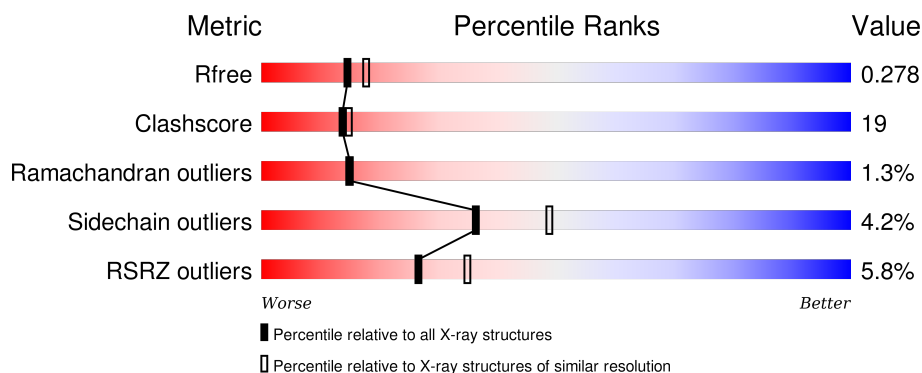
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	219	<div> <div>5%</div> <div> <div></div> <div>55%</div> <div>30%</div> <div>• • 12%</div> </div> </div>
1	B	219	<div> <div>4%</div> <div> <div></div> <div>60%</div> <div>26%</div> <div>• 11%</div> </div> </div>
1	C	219	<div> <div>4%</div> <div> <div></div> <div>64%</div> <div>23%</div> <div>• 10%</div> </div> </div>
1	D	219	<div> <div>8%</div> <div> <div></div> <div>54%</div> <div>32%</div> <div>• 12%</div> </div> </div>

## 2 Entry composition [i](#)

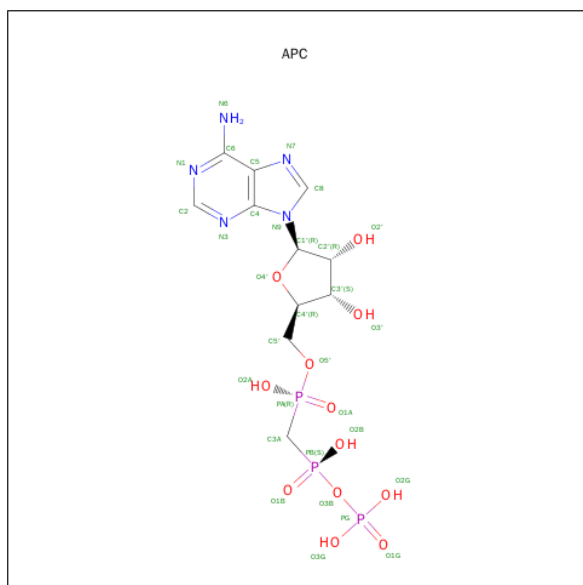
There are 5 unique types of molecules in this entry. The entry contains 6226 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called ADENYLATE CYCLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	193	Total	C	N	O	S	0	0	0
			1487	935	261	279	12			
1	B	194	Total	C	N	O	S	0	0	0
			1495	940	262	280	13			
1	C	198	Total	C	N	O	S	0	0	0
			1528	958	269	288	13			
1	D	193	Total	C	N	O	S	4	0	0
			1485	934	258	281	12			

- Molecule 2 is DIPHOSPHOMETHYLPHOSPHONIC ACID ADENOSYL ESTER (three-letter code: APC) (formula:  $C_{11}H_{18}N_5O_{12}P_3$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			31	11	5	12	3		
2	B	1	Total	C	N	O	P	0	0
			31	11	5	12	3		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	N	O	P	0	0
			31	11	5	12	3		
2	D	1	Total	C	N	O	P	0	0
			31	11	5	12	3		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		
3	A	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Ca	0	0
			1	1		
4	A	1	Total	Ca	0	0
			1	1		
4	D	1	Total	Ca	0	0
			1	1		
4	C	1	Total	Ca	0	0
			1	1		

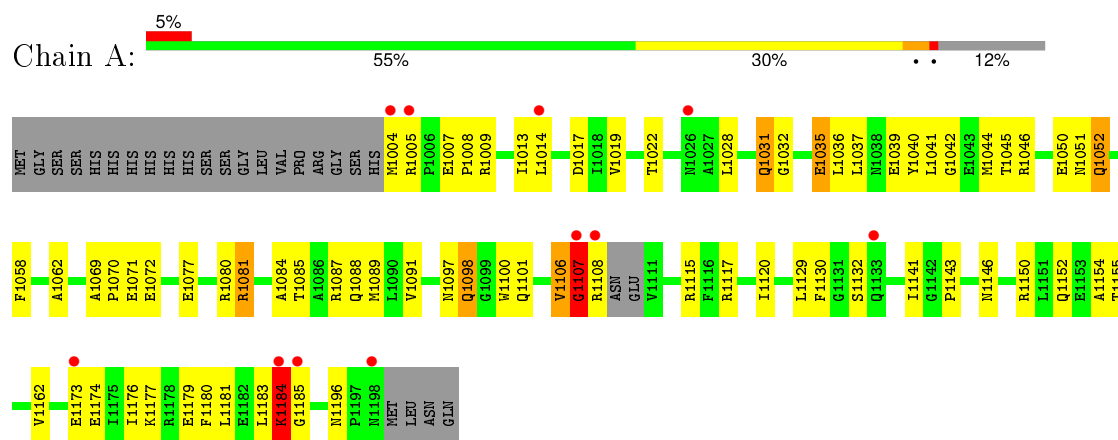
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	21	Total	O	0	0
			21	21		
5	B	36	Total	O	0	0
			36	36		
5	C	29	Total	O	0	0
			29	29		
5	D	13	Total	O	0	0
			13	13		

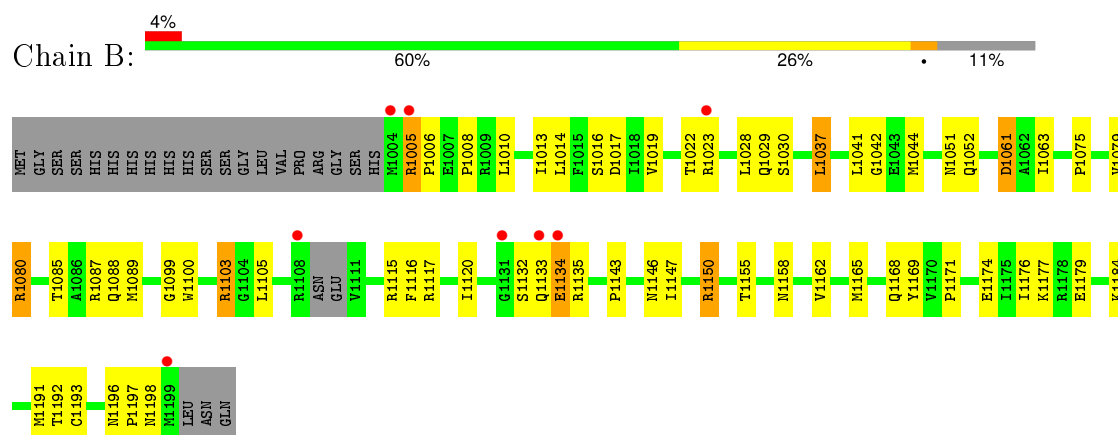
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

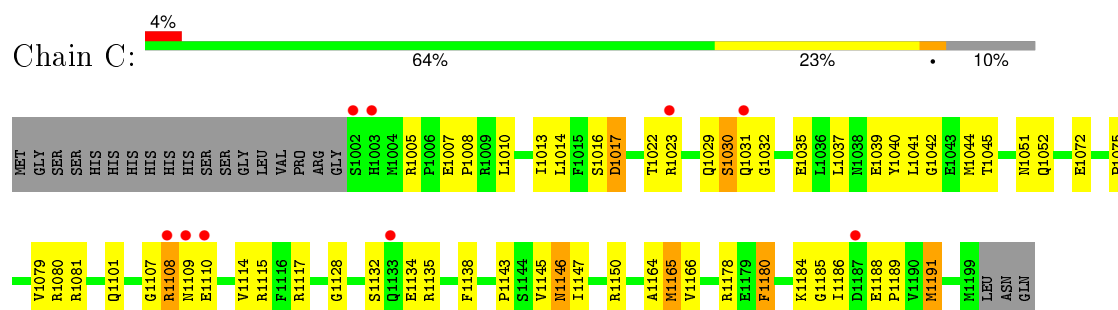
#### • Molecule 1: ADENYLATE CYCLASE



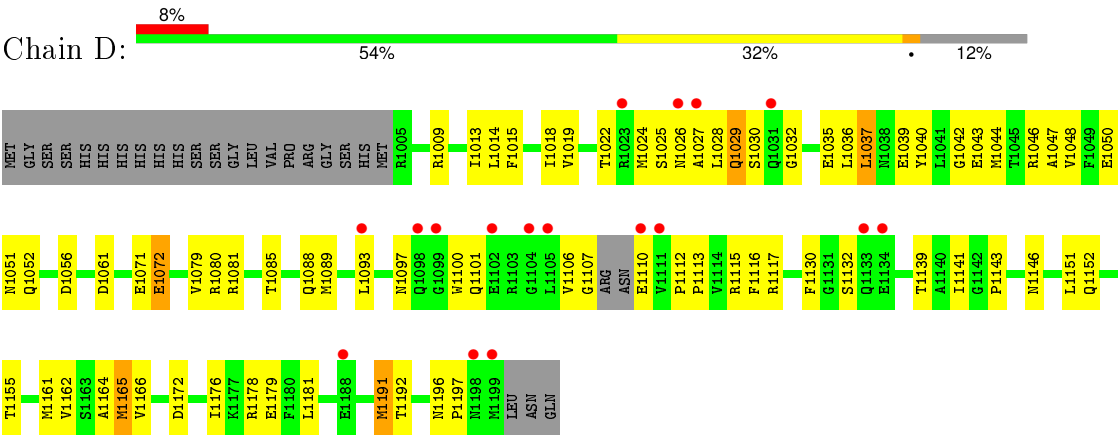
#### • Molecule 1: ADENYLATE CYCLASE



#### • Molecule 1: ADENYLATE CYCLASE



● Molecule 1: ADENYLATE CYCLASE



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	53.30 Å 71.19 Å 106.73 Å 90.00° 95.81° 90.00°	Depositor
Resolution (Å)	14.99 – 2.30 19.79 – 2.29	Depositor EDS
% Data completeness (in resolution range)	75.7 (14.99-2.30) 88.5 (19.79-2.29)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.50 (at 2.28 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.202 , 0.262 0.221 , 0.278	Depositor DCC
$R_{free}$ test set	2004 reflections (6.75%)	DCC
Wilson B-factor (Å <sup>2</sup> )	31.3	Xtriage
Anisotropy	0.481	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 40.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 33953 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	6226	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.25% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: APC, CA, MG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.36	0/1510	0.62	1/2041 (0.0%)
1	B	0.38	0/1518	0.62	0/2051
1	C	0.40	0/1553	0.61	1/2100 (0.0%)
1	D	0.33	0/1508	0.55	0/2039
All	All	0.37	0/6089	0.60	2/8231 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1107	GLY	N-CA-C	8.69	134.81	113.10
1	C	1017	ASP	N-CA-CB	-5.37	100.94	110.60

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1487	0	1498	76	0
1	B	1495	0	1507	60	0
1	C	1528	0	1532	47	0
1	D	1485	0	1491	58	0
2	A	31	0	14	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	31	0	14	0	0
2	C	31	0	14	0	0
2	D	31	0	14	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	21	0	0	2	0
5	B	36	0	0	2	0
5	C	29	0	0	2	0
5	D	13	0	0	1	0
All	All	6226	0	6084	226	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 19.

All (226) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1184:LYS:H	1:A:1184:LYS:HD3	1.07	1.18
1:A:1184:LYS:H	1:A:1184:LYS:CD	1.80	0.94
1:B:1134:GLU:HG3	1:B:1135:ARG:H	1.32	0.93
1:B:1005:ARG:HD2	1:B:1006:PRO:HD2	1.51	0.93
1:A:1031:GLN:NE2	1:A:1031:GLN:H	1.73	0.87
1:D:1018:ILE:HD12	1:D:1061:ASP:HB2	1.55	0.86
1:A:1150:ARG:NH1	1:A:1185:GLY:HA3	1.90	0.85
1:A:1184:LYS:N	1:A:1184:LYS:HD3	1.90	0.85
1:A:1154:ALA:HB2	1:A:1184:LYS:HE3	1.59	0.84
1:A:1008:PRO:O	1:A:1009:ARG:HG3	1.78	0.82
1:A:1101:GLN:OE1	1:A:1107:GLY:HA3	1.80	0.81
1:B:1150:ARG:HD3	1:B:1184:LYS:O	1.84	0.77
1:A:1005:ARG:HH12	1:A:1007:GLU:HB3	1.50	0.76
1:A:1150:ARG:HH11	1:A:1185:GLY:HA3	1.50	0.76
1:C:1134:GLU:HG2	1:C:1135:ARG:H	1.50	0.75
1:B:1042:GLY:HA2	1:C:1132:SER:HB3	1.69	0.75
1:A:1146:ASN:ND2	1:A:1150:ARG:HG2	2.03	0.73
1:D:1019:VAL:HG12	1:D:1115:ARG:O	1.87	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1031:GLN:HE21	1:A:1031:GLN:H	1.38	0.72
1:B:1080:ARG:HG2	1:B:1080:ARG:HH11	1.55	0.72
1:A:1042:GLY:HA2	1:D:1132:SER:HB3	1.72	0.71
1:A:1044:MET:HA	5:A:2004:HOH:O	1.90	0.71
1:D:1052:GLN:HB2	1:D:1081:ARG:HD3	1.73	0.70
1:B:1132:SER:HB3	1:C:1042:GLY:HA2	1.72	0.70
1:A:1132:SER:OG	1:D:1042:GLY:HA2	1.93	0.69
1:A:1037:LEU:HD12	1:D:1130:PHE:HD2	1.55	0.69
1:B:1134:GLU:HG3	1:B:1135:ARG:N	2.07	0.68
1:C:1014:LEU:C	1:C:1014:LEU:HD23	2.14	0.68
1:B:1019:VAL:HG11	1:B:1115:ARG:HB2	1.76	0.67
1:C:1180:PHE:CE1	1:C:1191:MET:HG2	2.30	0.67
1:C:1134:GLU:HG2	1:C:1135:ARG:N	2.10	0.66
1:B:1022:THR:HG23	1:C:1143:PRO:HG3	1.78	0.65
1:A:1005:ARG:NH1	1:A:1007:GLU:HB3	2.11	0.65
1:A:1141:ILE:HD13	1:D:1037:LEU:HD12	1.78	0.64
1:A:1085:THR:O	1:A:1089:MET:HG3	1.98	0.64
1:A:1087:ARG:O	1:A:1091:VAL:HG23	1.98	0.64
1:A:1146:ASN:HD21	1:A:1150:ARG:HG2	1.62	0.63
1:D:1093:LEU:HD21	1:D:1115:ARG:HA	1.81	0.63
1:D:1117:ARG:HB2	1:D:1152:GLN:HE21	1.63	0.63
1:D:1047:ALA:HA	1:D:1088:GLN:NE2	2.14	0.63
1:C:1165:MET:HB3	5:C:2025:HOH:O	1.98	0.63
1:B:1115:ARG:HB3	1:B:1158:ASN:ND2	2.13	0.62
1:A:1081:ARG:HH12	1:A:1084:ALA:HB3	1.63	0.62
1:D:1014:LEU:HD23	1:D:1014:LEU:C	2.20	0.62
1:D:1085:THR:O	1:D:1089:MET:HG3	1.99	0.61
1:A:1183:LEU:O	1:A:1185:GLY:N	2.34	0.61
1:A:1081:ARG:HA	1:A:1081:ARG:NH1	2.16	0.61
1:A:1019:VAL:HG12	1:A:1115:ARG:O	2.01	0.61
1:B:1117:ARG:HD3	1:B:1155:THR:O	2.00	0.60
1:B:1165:MET:O	1:B:1168:GLN:HG2	2.01	0.60
1:A:1177:LYS:HE3	1:A:1179:GLU:OE1	2.01	0.60
1:B:1087:ARG:HG2	1:B:1197:PRO:O	2.01	0.60
1:B:1099:GLY:O	1:B:1103:ARG:HB2	2.02	0.60
1:A:1117:ARG:HB2	1:A:1152:GLN:HE21	1.66	0.60
1:B:1085:THR:O	1:B:1089:MET:HG3	2.02	0.59
1:B:1116:PHE:H	1:B:1158:ASN:ND2	2.00	0.59
1:A:1035:GLU:O	1:A:1039:GLU:HG3	2.03	0.59
1:A:1037:LEU:HD12	1:D:1130:PHE:CD2	2.36	0.59
1:B:1005:ARG:HD2	1:B:1006:PRO:CD	2.31	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1018:ILE:HB	1:D:1061:ASP:OD2	2.02	0.59
1:D:1143:PRO:HD2	5:D:2007:HOH:O	2.03	0.58
1:B:1019:VAL:HG12	1:B:1115:ARG:O	2.03	0.58
1:A:1019:VAL:HG11	1:A:1115:ARG:HB2	1.86	0.58
1:D:1097:ASN:O	1:D:1101:GLN:HB2	2.02	0.58
1:C:1165:MET:SD	1:C:1165:MET:N	2.76	0.58
1:B:1010:LEU:CD2	1:B:1075:PRO:HG3	2.33	0.58
1:C:1035:GLU:O	1:C:1039:GLU:HG3	2.04	0.58
1:A:1041:LEU:HD22	1:A:1058:PHE:CE2	2.40	0.57
1:A:1040:TYR:O	1:A:1044:MET:HG2	2.05	0.57
1:D:1106:VAL:HG11	1:D:1113:PRO:HA	1.85	0.57
1:C:1052:GLN:HG3	1:C:1081:ARG:NE	2.20	0.57
1:C:1184:LYS:HG2	1:C:1185:GLY:N	2.18	0.57
1:C:1180:PHE:CD1	1:C:1191:MET:HG2	2.40	0.56
1:A:1007:GLU:HB2	1:A:1008:PRO:HD2	1.86	0.56
1:B:1150:ARG:HD2	5:B:2032:HOH:O	2.04	0.56
1:A:1031:GLN:HE21	1:A:1031:GLN:N	2.02	0.56
1:A:1143:PRO:HG3	1:D:1022:THR:HG23	1.87	0.56
1:B:1019:VAL:CG1	1:B:1115:ARG:HB2	2.35	0.56
1:D:1043:GLU:O	1:D:1046:ARG:HB3	2.06	0.56
1:B:1080:ARG:NH1	1:B:1080:ARG:HG2	2.21	0.55
1:B:1022:THR:CG2	1:C:1143:PRO:HG3	2.35	0.55
1:C:1134:GLU:N	1:C:1134:GLU:OE1	2.38	0.55
1:A:1081:ARG:HH12	1:A:1084:ALA:CB	2.20	0.55
1:B:1014:LEU:HD23	1:B:1014:LEU:C	2.27	0.54
1:A:1101:GLN:HE22	1:A:1108:ARG:N	2.05	0.54
1:A:1117:ARG:HD3	1:A:1155:THR:O	2.07	0.54
1:B:1030:SER:HB3	1:C:1008:PRO:HG3	1.90	0.54
1:C:1023:ARG:HH11	1:C:1023:ARG:HG3	1.73	0.54
1:A:1115:ARG:HH11	1:A:1115:ARG:HG3	1.73	0.54
1:D:1164:ALA:HB2	1:D:1191:MET:HG2	1.90	0.54
1:A:1019:VAL:CG1	1:A:1115:ARG:HB2	2.38	0.54
1:A:1028:LEU:O	1:A:1032:GLY:HA3	2.07	0.53
1:C:1052:GLN:HG3	1:C:1081:ARG:CZ	2.38	0.53
1:D:1026:ASN:O	1:D:1027:ALA:HB3	2.08	0.53
1:A:1004:MET:N	1:A:1129:LEU:HD12	2.24	0.53
1:B:1120:ILE:HB	1:B:1162:VAL:HG12	1.90	0.53
1:B:1017:ASP:O	1:B:1116:PHE:HB2	2.09	0.53
1:A:1150:ARG:HD2	1:A:1184:LYS:O	2.10	0.52
1:A:1184:LYS:N	1:A:1184:LYS:CD	2.57	0.52
1:A:1051:ASN:OD1	1:A:1081:ARG:NH1	2.43	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1128:GLY:HA2	1:C:1138:PHE:CE1	2.45	0.52
1:A:1141:ILE:HD13	1:D:1037:LEU:CD1	2.39	0.51
1:D:1040:TYR:O	1:D:1044:MET:HG2	2.10	0.51
1:C:1080:ARG:HG2	1:C:1080:ARG:HH11	1.75	0.51
1:B:1028:LEU:O	1:B:1029:GLN:HB2	2.11	0.51
1:D:1079:VAL:HG13	1:D:1166:VAL:HG13	1.91	0.51
1:A:1045:THR:HG21	1:A:1058:PHE:HZ	1.75	0.51
1:A:1081:ARG:NH1	1:A:1084:ALA:HB3	2.25	0.51
1:D:1014:LEU:HD23	1:D:1015:PHE:N	2.25	0.51
1:C:1013:ILE:HD13	1:C:1145:VAL:HA	1.92	0.51
1:B:1115:ARG:HA	1:B:1158:ASN:HD21	1.76	0.51
1:B:1005:ARG:HG3	1:B:1005:ARG:HH11	1.77	0.50
1:A:1017:ASP:OD1	1:A:1062:ALA:HB2	2.12	0.50
1:A:1071:GLU:OE2	1:A:1071:GLU:HA	2.12	0.50
1:A:1046:ARG:HD3	1:A:1050:GLU:OE2	2.12	0.50
1:C:1010:LEU:CD2	1:C:1075:PRO:HG3	2.42	0.50
1:B:1179:GLU:N	1:B:1191:MET:HE1	2.26	0.50
1:D:1035:GLU:O	1:D:1039:GLU:HG3	2.12	0.49
1:D:1036:LEU:HD23	1:D:1036:LEU:C	2.32	0.49
1:B:1051:ASN:C	1:B:1052:GLN:HG2	2.33	0.49
1:A:1098:GLN:O	1:A:1101:GLN:HB3	2.13	0.49
1:C:1164:ALA:HB2	1:C:1191:MET:HB3	1.95	0.49
1:C:1147:ILE:HG12	1:C:1186:ILE:HD13	1.95	0.49
1:B:1079:VAL:HG21	1:B:1169:TYR:HD1	1.78	0.49
1:C:1101:GLN:HE21	1:C:1107:GLY:HA3	1.78	0.49
1:A:1120:ILE:HB	1:A:1162:VAL:HG12	1.95	0.48
1:D:1151:LEU:HD22	1:D:1161:MET:HG2	1.95	0.48
1:B:1191:MET:HA	1:B:1191:MET:HE3	1.96	0.48
1:B:1191:MET:HE2	1:B:1192:THR:O	2.13	0.48
1:C:1178:ARG:NH1	1:C:1178:ARG:HG2	2.28	0.48
1:B:1088:GLN:HB2	5:B:2014:HOH:O	2.14	0.48
1:B:1150:ARG:CD	1:B:1184:LYS:O	2.59	0.48
1:A:1130:PHE:CZ	1:A:1141:ILE:HD12	2.49	0.47
1:D:1046:ARG:HG3	1:D:1050:GLU:OE2	2.14	0.47
1:B:1016:SER:HA	1:B:1117:ARG:O	2.15	0.47
1:B:1196:ASN:HD22	1:B:1197:PRO:HD2	1.78	0.47
1:A:1146:ASN:HD21	1:A:1150:ARG:CG	2.28	0.47
1:D:1107:GLY:O	1:D:1110:GLU:N	2.47	0.47
1:A:1176:ILE:CD1	1:A:1196:ASN:HA	2.45	0.47
1:C:1180:PHE:HE1	1:C:1191:MET:HG2	1.79	0.46
1:A:1176:ILE:HD11	1:A:1196:ASN:HA	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1146:ASN:HD21	1:C:1150:ARG:NH1	2.13	0.46
1:A:1150:ARG:HD3	1:A:1150:ARG:HA	1.79	0.46
1:B:1017:ASP:OD1	1:B:1061:ASP:OD2	2.34	0.46
1:D:1116:PHE:CD1	1:D:1116:PHE:C	2.89	0.46
1:A:1022:THR:HG22	2:A:1500:APC:O1B	2.15	0.46
1:B:1023:ARG:HH21	1:B:1023:ARG:HG3	1.80	0.46
1:C:1029:GLN:O	1:C:1032:GLY:N	2.42	0.46
1:D:1196:ASN:OD1	1:D:1197:PRO:HD2	2.16	0.45
1:D:1176:ILE:HD11	1:D:1196:ASN:HA	1.98	0.45
1:D:1018:ILE:HD11	1:D:1040:TYR:CD2	2.51	0.45
1:A:1084:ALA:O	1:A:1088:GLN:HG3	2.16	0.45
1:A:1044:MET:HE2	5:A:2004:HOH:O	2.15	0.45
1:C:1041:LEU:O	1:C:1045:THR:HG23	2.16	0.45
1:D:1009:ARG:NH2	1:D:1072:GLU:OE1	2.50	0.45
1:D:1028:LEU:C	1:D:1029:GLN:HG2	2.37	0.45
1:D:1115:ARG:HH11	1:D:1115:ARG:HG3	1.81	0.45
1:A:1179:GLU:OE2	1:A:1181:LEU:HD21	2.17	0.45
1:A:1014:LEU:C	1:A:1014:LEU:HD23	2.37	0.45
1:D:1164:ALA:HB3	1:D:1165:MET:CE	2.46	0.45
1:C:1010:LEU:HD23	1:C:1075:PRO:HG3	1.98	0.45
1:B:1019:VAL:O	1:B:1019:VAL:HG13	2.16	0.45
1:A:1130:PHE:CE2	1:A:1141:ILE:HD12	2.52	0.44
1:B:1100:TRP:CD1	1:B:1105:LEU:HD12	2.52	0.44
1:B:1177:LYS:O	1:B:1193:CYS:HA	2.17	0.44
1:B:1176:ILE:CD1	1:B:1196:ASN:HA	2.47	0.44
1:B:1146:ASN:O	1:B:1150:ARG:HG2	2.17	0.44
1:B:1115:ARG:CA	1:B:1158:ASN:HD21	2.31	0.44
1:D:1152:GLN:O	1:D:1155:THR:HG22	2.18	0.44
1:D:1044:MET:O	1:D:1048:VAL:HG23	2.18	0.44
2:A:1500:APC:H3A2	1:D:1146:ASN:ND2	2.33	0.44
1:D:1181:LEU:HD12	1:D:1192:THR:HG21	1.99	0.44
1:C:1040:TYR:O	1:C:1044:MET:HG2	2.17	0.44
1:D:1051:ASN:ND2	1:D:1085:THR:HA	2.33	0.44
1:C:1080:ARG:HG2	1:C:1080:ARG:NH1	2.33	0.44
1:A:1005:ARG:HH11	1:A:1005:ARG:HG2	1.83	0.43
1:A:1013:ILE:O	1:A:1120:ILE:HA	2.18	0.43
1:D:1022:THR:O	1:D:1025:SER:HB3	2.17	0.43
1:C:1178:ARG:HH11	1:C:1178:ARG:HG2	1.84	0.43
1:A:1072:GLU:HA	1:A:1072:GLU:OE1	2.17	0.43
1:B:1191:MET:HE2	1:B:1192:THR:N	2.34	0.43
1:B:1008:PRO:HG3	1:C:1030:SER:HB2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1112:PRO:HB2	1:D:1113:PRO:HD2	2.01	0.43
1:C:1016:SER:HA	1:C:1117:ARG:O	2.18	0.43
1:C:1007:GLU:HB3	5:C:2002:HOH:O	2.19	0.42
1:A:1077:GLU:OE2	1:A:1080:ARG:HD2	2.19	0.42
1:D:1080:ARG:HG2	1:D:1080:ARG:HH11	1.84	0.42
1:C:1029:GLN:O	1:C:1030:SER:C	2.58	0.42
1:B:1171:PRO:HB2	1:B:1174:GLU:HG3	2.02	0.42
1:A:1100:TRP:HB3	1:A:1106:VAL:HG22	2.02	0.42
1:A:1037:LEU:O	1:A:1041:LEU:HG	2.19	0.42
1:D:1056:ASP:OD2	1:D:1139:THR:HA	2.19	0.42
1:B:1132:SER:O	1:B:1134:GLU:N	2.52	0.42
1:C:1014:LEU:C	1:C:1014:LEU:CD2	2.85	0.42
1:C:1108:ARG:HG3	1:C:1109:ASN:N	2.35	0.42
1:D:1101:GLN:NE2	1:D:1106:VAL:O	2.52	0.42
1:C:1051:ASN:C	1:C:1052:GLN:HG2	2.40	0.42
1:D:1028:LEU:O	1:D:1032:GLY:HA3	2.20	0.42
1:C:1188:GLU:HA	1:C:1189:PRO:HD3	1.95	0.42
1:C:1017:ASP:C	1:C:1017:ASP:OD2	2.57	0.42
1:B:1143:PRO:O	1:B:1147:ILE:HG13	2.20	0.42
1:D:1024:MET:O	1:D:1028:LEU:HB2	2.20	0.41
2:A:1500:APC:HN61	1:D:1141:ILE:HD13	1.85	0.41
1:A:1069:ALA:HA	1:A:1070:PRO:C	2.40	0.41
1:B:1116:PHE:H	1:B:1158:ASN:HD22	1.68	0.41
1:A:1097:ASN:O	1:A:1101:GLN:HB2	2.20	0.41
1:B:1143:PRO:HB3	1:C:1022:THR:HG23	2.02	0.41
1:D:1100:TRP:HB3	1:D:1106:VAL:HG23	2.02	0.41
1:B:1171:PRO:HD2	1:B:1174:GLU:OE1	2.20	0.41
1:A:1051:ASN:C	1:A:1052:GLN:HG2	2.41	0.41
1:B:1013:ILE:N	1:B:1013:ILE:HD12	2.36	0.41
1:B:1008:PRO:HG3	1:C:1030:SER:CB	2.50	0.41
1:D:1071:GLU:OE2	1:D:1071:GLU:HA	2.21	0.41
1:C:1114:VAL:HG12	1:C:1115:ARG:N	2.36	0.41
1:D:1176:ILE:CD1	1:D:1196:ASN:HA	2.51	0.41
1:A:1036:LEU:HD23	1:A:1036:LEU:C	2.41	0.41
1:B:1044:MET:CB	1:B:1063:ILE:CD1	2.98	0.41
1:A:1173:GLU:HG2	1:A:1174:GLU:N	2.35	0.41
1:A:1005:ARG:HG2	1:A:1005:ARG:NH1	2.36	0.40
1:D:1162:VAL:HB	1:D:1166:VAL:HB	2.03	0.40
1:D:1013:ILE:HD12	1:D:1013:ILE:N	2.35	0.40
1:C:1079:VAL:HG13	1:C:1166:VAL:HG13	2.03	0.40
1:D:1106:VAL:HG13	1:D:1112:PRO:O	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1044:MET:HB3	1:B:1063:ILE:CD1	2.51	0.40
1:D:1172:ASP:OD1	1:D:1178:ARG:NH2	2.54	0.40
1:B:1037:LEU:HD22	1:B:1041:LEU:CD1	2.51	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	189/219 (86%)	177 (94%)	9 (5%)	3 (2%)	12	11
1	B	190/219 (87%)	179 (94%)	7 (4%)	4 (2%)	9	7
1	C	196/219 (90%)	189 (96%)	6 (3%)	1 (0%)	34	41
1	D	189/219 (86%)	173 (92%)	14 (7%)	2 (1%)	17	18
All	All	764/876 (87%)	718 (94%)	36 (5%)	10 (1%)	15	15

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1184	LYS
1	B	1134	GLU
1	C	1030	SER
1	D	1029	GLN
1	A	1107	GLY
1	D	1030	SER
1	B	1103	ARG
1	B	1198	ASN
1	B	1133	GLN
1	A	1106	VAL

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	160/183 (87%)	153 (96%)	7 (4%)	35	46
1	B	161/183 (88%)	156 (97%)	5 (3%)	47	64
1	C	165/183 (90%)	155 (94%)	10 (6%)	23	30
1	D	160/183 (87%)	155 (97%)	5 (3%)	47	64
All	All	646/732 (88%)	619 (96%)	27 (4%)	36	49

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1031	GLN
1	A	1035	GLU
1	A	1052	GLN
1	A	1081	ARG
1	A	1098	GLN
1	A	1180	PHE
1	A	1184	LYS
1	B	1005	ARG
1	B	1037	LEU
1	B	1061	ASP
1	B	1080	ARG
1	B	1150	ARG
1	C	1005	ARG
1	C	1031	GLN
1	C	1037	LEU
1	C	1072	GLU
1	C	1108	ARG
1	C	1110	GLU
1	C	1146	ASN
1	C	1165	MET
1	C	1180	PHE
1	C	1191	MET
1	D	1037	LEU
1	D	1072	GLU

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Mol	Chain	Res	Type
1	D	1165	MET
1	D	1179	GLU
1	D	1191	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1031	GLN
1	A	1052	GLN
1	A	1122	GLN
1	A	1146	ASN
1	A	1152	GLN
1	A	1158	ASN
1	B	1052	GLN
1	B	1122	GLN
1	B	1152	GLN
1	B	1158	ASN
1	B	1196	ASN
1	C	1098	GLN
1	C	1101	GLN
1	C	1146	ASN
1	C	1152	GLN
1	C	1158	ASN
1	D	1088	GLN
1	D	1101	GLN
1	D	1152	GLN
1	D	1158	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

Of 12 ligands modelled in this entry, 8 are monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z  > 2$	Counts	RMSZ	# $ Z  > 2$
2	APC	A	1500	3,4	25,33,33	1.60	5 (20%)	30,52,52	1.19	4 (13%)
2	APC	B	1500	3,4	25,33,33	1.79	5 (20%)	30,52,52	1.13	3 (10%)
2	APC	C	1500	3,4	25,33,33	1.58	4 (16%)	30,52,52	1.27	4 (13%)
2	APC	D	1500	3,4	25,33,33	1.50	5 (20%)	30,52,52	1.15	3 (10%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	APC	A	1500	3,4	-	0/15/38/38	0/3/3/3
2	APC	B	1500	3,4	-	0/15/38/38	0/3/3/3
2	APC	C	1500	3,4	-	0/15/38/38	0/3/3/3
2	APC	D	1500	3,4	-	0/15/38/38	0/3/3/3

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1500	APC	C8-N7	-2.20	1.30	1.34
2	A	1500	APC	C8-N7	-2.15	1.30	1.34
2	D	1500	APC	C8-N7	-2.06	1.30	1.34
2	C	1500	APC	C8-N7	-2.00	1.30	1.34
2	C	1500	APC	C2-N1	2.18	1.38	1.33
2	A	1500	APC	C2-N1	2.23	1.38	1.33
2	B	1500	APC	C2-N1	2.35	1.38	1.33
2	D	1500	APC	C2-N1	2.54	1.38	1.33
2	B	1500	APC	PB-O3B	2.65	1.61	1.58
2	D	1500	APC	PB-O3B	2.79	1.61	1.58
2	A	1500	APC	PB-O3B	3.23	1.62	1.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1500	APC	O4'-C1'	3.36	1.45	1.41
2	D	1500	APC	PA-O5'	3.48	1.61	1.57
2	C	1500	APC	PA-O5'	3.58	1.61	1.57
2	D	1500	APC	O4'-C1'	3.98	1.46	1.41
2	B	1500	APC	O4'-C1'	4.00	1.46	1.41
2	A	1500	APC	PA-O5'	4.68	1.62	1.57
2	C	1500	APC	O4'-C1'	5.04	1.47	1.41
2	B	1500	APC	PA-O5'	5.90	1.64	1.57

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1500	APC	PG-O3B-PB	-2.98	122.68	132.67
2	D	1500	APC	PG-O3B-PB	-2.88	122.99	132.67
2	C	1500	APC	PG-O3B-PB	-2.83	123.17	132.67
2	C	1500	APC	C2'-C1'-N9	-2.80	110.02	114.29
2	C	1500	APC	O5'-PA-O1A	-2.63	107.00	113.98
2	B	1500	APC	PG-O3B-PB	-2.57	124.06	132.67
2	D	1500	APC	O5'-PA-O1A	-2.33	107.79	113.98
2	B	1500	APC	C2'-C1'-N9	-2.16	111.00	114.29
2	A	1500	APC	C1'-N9-C4	-2.13	123.73	126.94
2	A	1500	APC	O5'-PA-O1A	-2.04	108.54	113.98
2	D	1500	APC	C4-C5-N7	2.27	111.57	109.48
2	A	1500	APC	C4-C5-N7	2.40	111.69	109.48
2	B	1500	APC	C4-C5-N7	2.53	111.80	109.48
2	C	1500	APC	C4-C5-N7	2.53	111.81	109.48

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1500	APC	3	0

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å <sup>2</sup> )	Q<0.9
1	A	193/219 (88%)	0.41	11 (5%)	27 36	23, 39, 65, 92	0
1	B	194/219 (88%)	0.09	8 (4%)	41 50	13, 29, 60, 77	0
1	C	198/219 (90%)	0.16	9 (4%)	37 46	14, 31, 62, 67	0
1	D	193/219 (88%)	0.47	17 (8%)	12 18	20, 43, 81, 89	1 (0%)
All	All	778/876 (88%)	0.28	45 (5%)	26 35	13, 35, 67, 92	1 (0%)

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1004	MET	5.6
1	B	1004	MET	5.5
1	A	1108	ARG	5.3
1	B	1133	GLN	4.4
1	B	1108	ARG	4.2
1	D	1111	VAL	4.1
1	A	1133	GLN	3.9
1	B	1199	MET	3.9
1	D	1027	ALA	3.7
1	C	1133	GLN	3.6
1	D	1105	LEU	3.6
1	A	1198	ASN	3.6
1	D	1098	GLN	3.6
1	C	1187	ASP	3.5
1	A	1107	GLY	3.4
1	A	1005	ARG	3.2
1	D	1104	GLY	3.2
1	D	1023	ARG	3.2
1	D	1110	GLU	3.2
1	D	1133	GLN	3.1
1	C	1110	GLU	3.1

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Mol	Chain	Res	Type	RSRZ
1	D	1188	GLU	2.9
1	B	1131	GLY	2.8
1	C	1109	ASN	2.8
1	D	1199	MET	2.7
1	D	1102	GLU	2.6
1	A	1173	GLU	2.6
1	B	1134	GLU	2.5
1	A	1026	ASN	2.5
1	C	1002	SER	2.4
1	C	1023	ARG	2.3
1	C	1003	HIS	2.3
1	C	1031	GLN	2.3
1	B	1005	ARG	2.3
1	D	1026	ASN	2.2
1	D	1099	GLY	2.2
1	D	1134	GLU	2.2
1	D	1198	ASN	2.2
1	A	1184	LYS	2.1
1	B	1023	ARG	2.1
1	A	1185	GLY	2.1
1	C	1108	ARG	2.1
1	A	1014	LEU	2.0
1	D	1031	GLN	2.0
1	D	1093	LEU	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	APC	A	1500	31/31	0.95	0.13	-0.54	39,41,44,46	0
2	APC	B	1500	31/31	0.96	0.10	-0.74	20,25,34,35	0
2	APC	D	1500	31/31	0.95	0.11	-0.77	34,38,43,47	0
2	APC	C	1500	31/31	0.98	0.09	-0.87	16,25,30,31	0
4	CA	C	2100	1/1	0.98	0.06	-2.30	24,24,24,24	0
4	CA	B	2100	1/1	0.96	0.05	-3.29	29,29,29,29	0
4	CA	D	2100	1/1	0.99	0.02	-6.33	32,32,32,32	0
3	MG	A	2000	1/1	0.92	0.08	-	42,42,42,42	0
3	MG	C	2000	1/1	0.95	0.08	-	18,18,18,18	0
3	MG	D	2000	1/1	0.92	0.07	-	40,40,40,40	0
3	MG	B	2000	1/1	0.88	0.09	-	27,27,27,27	0
4	CA	A	2100	1/1	0.98	0.03	-	32,32,32,32	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.